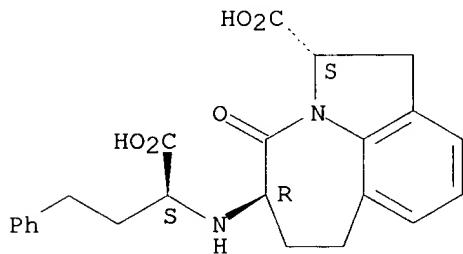


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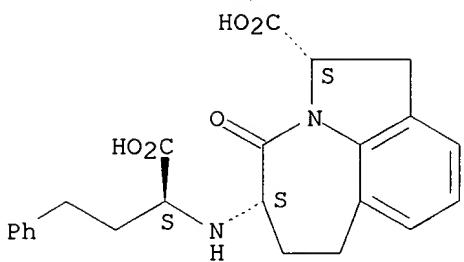
L23 ANSWER 90 OF 167 CAPLUS COPYRIGHT 2002 ACS  
AN 1995:248136 CAPLUS  
DN 122:133800  
TI Practical syntheses of a novel tricyclic dipeptide mimetic based on a [6H]-azepinoindoline nucleus: application to angiotensin-converting enzyme inhibition  
AU De Lombaert, Stephane; Blanchard, Louis; Stamford, Lisa B.; Sperbeck, Donald M.; Grim, Michael D.; Jenson, Todd M.; Rodriguez, Herman R.  
CS Res. Dep., CIBA-GEIGY Corp., Summit, NJ, 07901, USA  
SO Tetrahedron Lett. (1994), 35(41), 7513-16  
CODEN: TELEAY; ISSN: 0040-4039  
DT Journal  
LA English  
OS CASREACT 122:133800  
AB Two stereocontrolled synthetic approaches towards 5(S)-amino-1,2,4,5,6,7-hexahydro-4-oxo-azepino[3,2,1-h]indole-2(S)-carboxylic acid, based on intramol. Friedel-Crafts acylations, are reported. This conformationally restricted tricyclic dipeptidomimetic has been converted to its 5-[1(S)-carboxy-3-phenylpropyl]amino analog, a potent and orally active inhibitor of angiotensin-converting enzyme (ACE).  
IT 160726-16-9P 160798-20-9P  
RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)  
(prepn. of a tricyclic dipeptide mimetic as ACE inhibitor)  
RN 160726-16-9 CAPLUS  
CN Azepino[3,2,1-h]indole-2-carboxylic acid, 5-[(1-carboxy-3-phenylpropyl)amino]-1,2,4,5,6,7-hexahydro-4-oxo-, [2S-[2.alpha.,5.beta.(R\*)]]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 160798-20-9 CAPLUS  
CN Azepino[3,2,1-h]indole-2-carboxylic acid, 5-[(1-carboxy-3-phenylpropyl)amino]-1,2,4,5,6,7-hexahydro-4-oxo-, [2S-[2.alpha.,5.alpha.(R\*)]]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



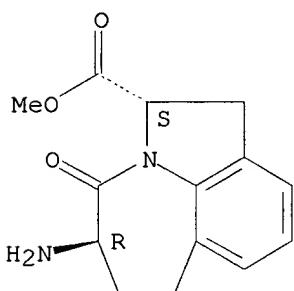
IT 160865-33-8P

RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
(prepn. of a tricyclic dipeptide mimetic as ACE inhibitor)

RN 160865-33-8 CAPLUS

CN Azepino[3,2,1-hi]indole-2-carboxylic acid,  
5-amino-1,2,4,5,6,7-hexahydro-4-  
oxo-, methyl ester, monohydrobromide, (2S-trans)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● HBr

IT 160726-07-8P 160726-08-9P 160726-09-0P

160798-17-4P 160798-18-5P 160798-19-6P

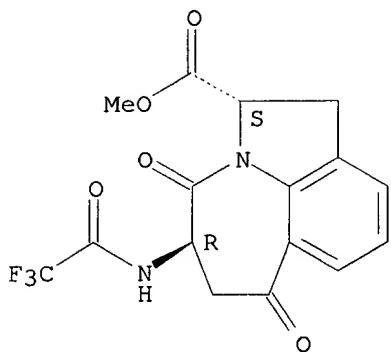
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation)  
(prepn. of a tricyclic dipeptide mimetic as ACE inhibitor)

RN 160726-07-8 CAPLUS

CN Azepino[3,2,1-hi]indole-2-carboxylic acid,  
1,2,4,5,6,7-hexahydro-4,7-dioxo-  
5-[ (trifluoroacetyl)amino]-, methyl ester, (2S-trans)- (9CI) (CA INDEX  
NAME)

Absolute stereochemistry.

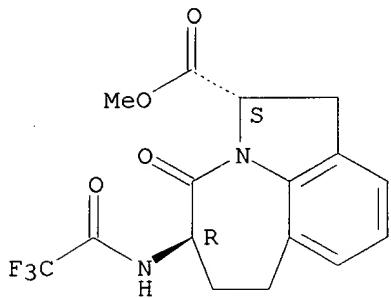
09/882,777



RN 160726-08-9 CAPLUS

CN Azepino[3,2,1-hi]indole-2-carboxylic acid, 1,2,4,5,6,7-hexahydro-4-oxo-5-[(trifluoroacetyl)amino]-, methyl ester, (2S-trans)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

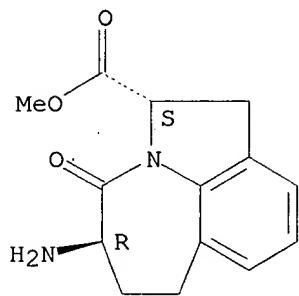


RN 160726-09-0 CAPLUS

CN Azepino[3,2,1-hi]indole-2-carboxylic acid,  
5-amino-1,2,4,5,6,7-hexahydro-4-  
oxo-, methyl ester, monohydrochloride, (2S-trans)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

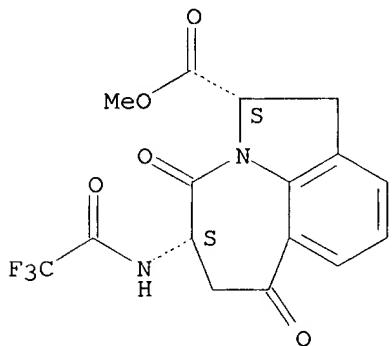
09/882,777



● HCl

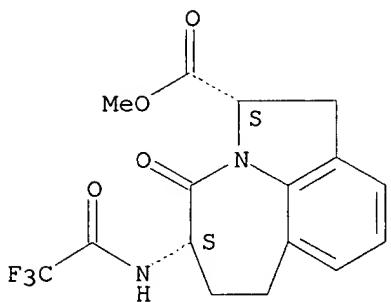
RN 160798-17-4 CAPLUS  
CN Azepino[3,2,1-hi]indole-2-carboxylic acid,  
1,2,4,5,6,7-hexahydro-4,7-dioxo-  
5-[(trifluoroacetyl)amino]-, methyl ester, (2S-cis)- (9CI) (CA INDEX  
NAME)

Absolute stereochemistry.



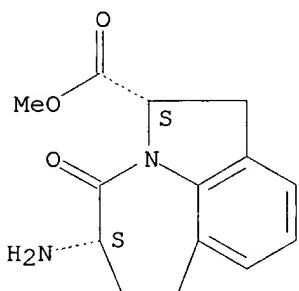
RN 160798-18-5 CAPLUS  
CN Azepino[3,2,1-hi]indole-2-carboxylic acid, 1,2,4,5,6,7-hexahydro-4-oxo-5-  
[(trifluoroacetyl)amino]-, methyl ester, (2S-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



RN 160798-19-6 CAPLUS  
CN Azepino[3,2,1-hi]indole-2-carboxylic acid,  
5-amino-1,2,4,5,6,7-hexahydro-4-  
oxo-, methyl ester, monohydrochloride, (2S-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

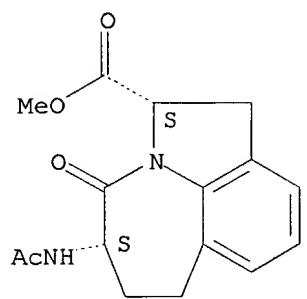


● HCl

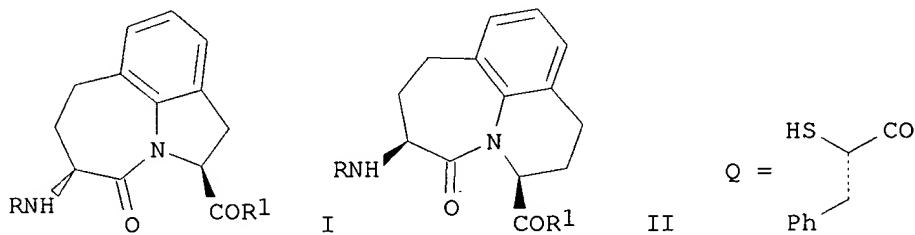
IT 160726-15-8P  
RL: SPN (Synthetic preparation); PREP (Preparation)  
(prep. of a tricyclic dipeptide mimetic as ACE inhibitor)  
RN 160726-15-8 CAPLUS  
CN Azepino[3,2,1-hi]indole-2-carboxylic acid, 5-(acetylamino)-1,2,4,5,6,7-  
hexahydro-4-oxo-, methyl ester, (2S-cis)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

09/882,777



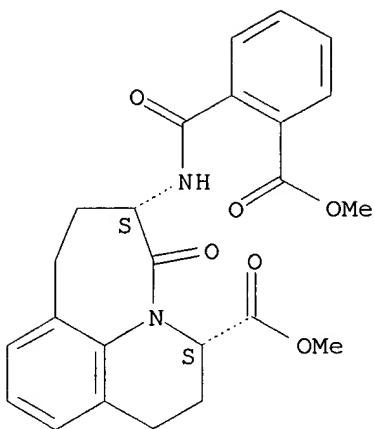
L23 ANSWER 85 OF 167 CAPLUS COPYRIGHT 2002 ACS  
 AN 1995:439878 CAPLUS  
 DN 123:112686  
 TI Synthesis of benzo-fused, 7,5- and 7,6-fused azepinones as conformationally restricted dipeptide mimetics  
 AU Robl, Jeffrey A.; Karanewsky, Donald S.; Asaad, Magdi M.  
 CS Bristol-Myers Squibb Pharmaceutical Res. Institute, Princeton, NJ, 08453-4000, USA  
 SO Tetrahedron Letters (1995), 36(10), 1593-6  
 CODEN: TELEAY; ISSN: 0040-4039  
 PB Elsevier  
 DT Journal  
 LA English  
 OS CASREACT 123:112686  
 GI



AB Methodol. for the generation of novel conformationally restricted dipeptide mimetics I and II (R = H, R<sup>1</sup> = OEt) has been developed. The key step involved intramol. addn. of an oxonium ion to the proximal indoline/tetrahydroquinoline arom. ring. A dramatic difference in reactivity was obsd. in the formation of the 7,5-vs. the 7,6-fused azepinone nuclei. Application of these mimetics in the synthesis of dual-acting angiotensin converting enzyme (ACE)/neutral endopeptidase (NEP) inhibitors I and II (R = Q, R<sup>1</sup> = OH) is described.

IT **165748-16-3P**  
 RL: BYP (Byproduct); PREP (Preparation)  
 (prep. of aminoazepinoindolecarboxylate  
 aminopyridobenzazepinecarboxyl  
 ate derivs. as conformationally constrained dipeptide mimics)  
 RN 165748-16-3 CAPLUS  
 CN 1H-Pyrido[3,2,1-jk][1]benzazepine-3-carboxylic acid,  
 2,3,5,6,7,8-hexahydro-  
 6-[(2-(methoxycarbonyl)benzoyl)amino]-5-oxo-, methyl ester, (3S-cis)-  
 (9CI) (CA INDEX NAME)

Relative stereochemistry.



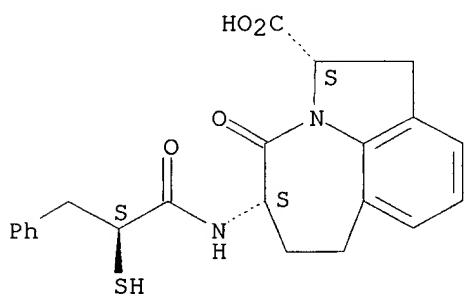
IT 157521-10-3P 157521-15-8P

RL: PNU (Preparation, unclassified); PREP (Preparation)  
(prepn. of aminoazepinoindolecarboxylate  
aminopyridobenzazepinecarboxyl  
ate derivs. as conformationally constrained dipeptide mimics)

RN 157521-10-3 CAPLUS

CN Azepino[3,2,1-hi]indole-2-carboxylic acid,  
1,2,4,5,6,7-hexahydro-5-[(2S)-  
2-mercaptop-1-oxo-3-phenylpropyl]amino]-4-oxo-, (2S,5S)- (9CI) (CA INDEX  
NAME)

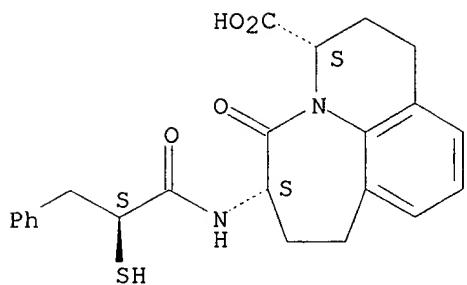
Absolute stereochemistry. Rotation (+).



RN 157521-15-8 CAPLUS

CN 1H-Pyrido[3,2,1-jk][1]benzazepine-3-carboxylic acid,  
2,3,5,6,7,8-hexahydro-  
6-(2-mercaptop-1-oxo-3-phenylpropyl)-5-oxo-, [3S-[3.alpha.,6.alpha.(R\*)]]-  
(9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



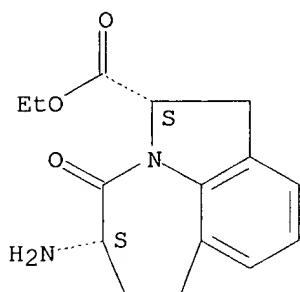
IT 157521-25-0P 166020-24-2P

RL: SPN (Synthetic preparation); PREP (Preparation)  
 (prepn. of aminoazepinoindolecarboxylate  
 aminopyridobenzazepinecarboxyl  
 ate derivs. as conformationally constrained dipeptide mimics)

RN 157521-25-0 CAPLUS

CN Azepino[3,2,1-h]indole-2-carboxylic acid,  
 5-amino-1,2,4,5,6,7-hexahydro-4-  
 oxo-, ethyl ester, (2S,5S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



RN 166020-24-2 CAPLUS